

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX121322\
 Data File : VX033396.D
 Acq On : 13 Dec 2022 16:40
 Operator : JC/MD
 Sample : N6001-11
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-11B

Quant Time: Dec 14 05:39:35 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120722W.M
 Quant Title : SW846 8260
 QLast Update : Wed Dec 07 12:43:54 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Pentafluorobenzene	5.556	168	123485	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	196145	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	176382	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	79947	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	38000	47.968	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	95.940%
35) Dibromofluoromethane	5.391	113	42348	48.002	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	96.000%
50) Toluene-d8	8.647	98	90729	49.161	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	98.320%
62) 4-Bromofluorobenzene	11.079	95	68677	48.668	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	97.340%
Target Compounds						
					Qvalue	
40) Benzene	6.043	78	13882	2.768	ug/l	98
67) Ethyl Benzene	10.195	91	3194	0.513	ug/l #	87
68) m/p-Xylenes	10.305	106	2277	0.941	ug/l	93
69) o-Xylene	10.646	106	603	0.252	ug/l	74

(#) = qualifier out of range (m) = manual integration (+) = signals summed

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX121322\
Data File : VX033396.D
Acq On : 13 Dec 2022 16:40
Operator : JC/MD
Sample : N6001-11
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-11B

Quant Time: Dec 14 05:39:35 2022
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120722W.M
Quant Title : SW846 8260
QLast Update : Wed Dec 07 12:43:54 2022
Response via : Initial Calibration

